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TITLE: Benzo-y-pyrones. Part XIV. Reaction of

C-substituted 2-phenyl-4H-1-benzopyran-4-ones with

hydroxylamine

AUTHOR(S): Basinski, Wlodzimierz

CORPORATE SOURCE: Fac. Pharm., Sch. Med., Lodz, 90151, Pol.

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GΙ

$$R^{1}$$
 R^{2}
 R^{3}
 R^{1}
 R^{2}
 R^{2}
 R^{3}
 R^{2}
 R^{3}
 R^{3}
 R^{1}
 R^{2}
 R^{3}
 R^{3}
 R^{1}
 R^{2}
 R^{3}
 R^{3}

AB The reaction of flavones I (R, R1 = H, Me; R2 = H, Me, Br; R3 = H, MeO) with hydroxylamine in anhydrous pyridine was investigated. The oximes II and isoxazoles III were the products. It was determined that the ratio of II to III is dependent on the nature of substituent and its position in the flavone skeleton. It is postulated that the flavone is an ambient electrophile and that the reaction course is characteristic for this class of compds.

IT 115663-23-5P 140885-79-6P

RL: SPN (Synthetic preparation); PREP (Preparation) (preparation, acetylation, IR, NMR, and mass spectrum of)

RN 115663-23-5 CAPLUS

CN 4H-1-Benzopyran-4-one, 2-(4-methoxyphenyl)-6-methyl-, oxime (CA INDEX NAME)

RN 140885-79-6 CAPLUS

CN 4H-1-Benzopyran-4-one, 8-methyl-2-phenyl-, oxime (CA INDEX NAME)